

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
 Data File : BP009412.D
 Acq On : 15 Mar 2022 23:01
 Operator : CG/JU
 Sample : N1894-02 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP2(0-5)

Quant Time: Mar 16 01:59:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian
 Giraldo

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							03/16/2022
1) 1,4-Dichlorobenzene-d4	7.810	152	319958	20.000	ng	-0.01	Supervised By :Jagrut Upadhyay
21) Naphthalene-d8	10.604	136	1280229	20.000	ng	-0.01	
39) Acenaphthene-d10	14.451	164	858297	20.000	ng	-0.01	
64) Phenanthrene-d10	17.216	188	1819596	20.000	ng	0.00	
76) Chrysene-d12	21.327	240	1767325	20.000	ng	0.00	
86) Perylene-d12	23.745	264	1978938	20.000	ng	0.01	03/16/2022
System Monitoring Compounds							
5) 2-Fluorophenol	5.422	112	824506	37.615	ng	0.01	
7) Phenol-d6	7.005	99	1014544	36.422	ng	0.02	
23) Nitrobenzene-d5	8.963	82	657936	28.306	ng	-0.01	
42) 2,4,6-Tribromophenol	15.951	330	608000	58.665	ng	0.00	
45) 2-Fluorobiphenyl	13.075	172	1675955	26.853	ng	-0.01	
79) Terphenyl-d14	19.792	244	2783240	29.719	ng	0.00	
Target Compounds							Qvalue
49) Acenaphthylene	14.175	152	532509	6.493	ng		99
71) Phenanthrene	17.257	178	1395708	13.347	ng		99
72) Anthracene	17.351	178	536903	5.158	ng		99
73) Carbazole	17.628	167	277336	2.924	ng		99
75) Fluoranthene	19.239	202	3493860	28.946	ng		98
78) Pyrene	19.592	202	3049110	25.553	ng		100
81) Benzo(a)anthracene	21.310	228	1872932	15.635	ng		97
83) Chrysene	21.363	228	1710458	14.707	ng		98
84) Bis(2-ethylhexyl)phtha...	21.245	149	346766	5.034	ng	#	97
87) Indeno(1,2,3-cd)pyrene	26.251	276	1546380	10.058	ng	#	91
88) Benzo(b)fluoranthene	23.010	252	2699400m	20.203	ng		
89) Benzo(k)fluoranthene	23.051	252	892057m	6.988	ng		
90) Benzo(a)pyrene	23.633	252	1805287	14.245	ng		96
91) Dibenzo(a,h)anthracene	26.256	278	415381	3.186	ng		94
92) Benzo(g,h,i)perylene	27.021	276	1488668	11.558	ng		95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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