

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072320\
 Data File : BP002842.D
 Acq On : 23 Jul 2020 16:11
 Operator : CG/JU
 Sample : L3360-01RX
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP-1ARX

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:13:13 PM

Quant Time: Jul 24 02:18:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	62986	20.00	ng	-0.01
21) Naphthalene-d8	10.31	136	274197	20.00	ng	-0.02
39) Acenaphthene-d10	14.20	164	175983	20.00	ng	-0.01
64) Phenanthrene-d10	16.96	188	416574	20.00	ng	-0.01
76) Chrysene-d12	21.07	240	516122	20.00	ng	0.00
86) Perylene-d12	23.26	264	594666	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	161853	43.36	ng	-0.01
7) Phenol-d6	6.75	99	452297	84.89	ng	-0.01
23) Nitrobenzene-d5	8.69	82	326114	63.61	ng	-0.02
42) 2,4,6-Tribromophenol	15.70	330	57047	31.05	ng	0.00
45) 2-Fluorobiphenyl	12.80	172	662622	58.25	ng	-0.02
79) Terphenyl-d14	19.55	244	1088413	48.35	ng	-0.01
Target Compounds						
50) Dimethylphthalate	13.66	163	125198	9.328	ng	99
71) Phenanthrene	17.00	178	508638	22.551	ng	99
72) Anthracene	17.09	178	139127	6.286	ng	97
73) Carbazole	17.37	167	96224	4.430	ng	100
75) Fluoranthene	18.99	202	1565316	59.894	ng	98
78) Pyrene	19.35	202	1428225	40.079	ng	99
81) Benzo(a)anthracene	21.06	228	956929	28.642	ng	100
83) Chrysene	21.10	228	953619m	29.986	ng	
87) Indeno(1,2,3-cd)pyrene	25.47	276	513438	11.946	ng	96
88) Benzo(b)fluoranthene	22.61	252	1228124m	33.260	ng	
89) Benzo(k)fluoranthene	22.65	252	364198m	10.013	ng	
90) Benzo(a)pyrene	23.17	252	835443	23.876	ng	99
91) Dibenzo(a,h)anthracene	25.47	278	141935	4.095	ng	# 79
92) Benzo(g,h,i)perylene	26.14	276	504687	14.361	ng	97

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

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