

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081720\
 Data File : BP003009.D
 Acq On : 17 Aug 2020 18:53
 Operator : CG/JU
 Sample : L3691-10
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 S711-N-SO-0-0.5-081320-FD

Manual Integrations
 APPROVED

mohammad
 8/19/2020 2:20:51 PM

Quant Time: Aug 18 04:26:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 14:07:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	345510	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1299376	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	833057	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	1569099	20.00	ng	0.00
76) Chrysene-d12	21.17	240	1837422	20.00	ng	0.00
86) Perylene-d12	23.42	264	2020909	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1834115	81.25	ng	0.00
7) Phenol-d6	6.86	99	2196714	70.09	ng	0.00
23) Nitrobenzene-d5	8.83	82	1248660	52.05	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	835885	80.73	ng	0.00
45) 2-Fluorobiphenyl	12.95	172	3076616	50.46	ng	0.00
79) Terphenyl-d14	19.65	244	4082452	44.48	ng	0.00
Target Compounds						
10) Phenol	6.89	94	85619	2.712	ng	97
32) Benzoic acid	9.74	122	4527m	8.001	ng	
50) Dimethylphthalate	13.78	163	739787	11.354	ng	100
71) Phenanthrene	17.12	178	425569	4.625	ng	97
75) Fluoranthene	19.09	202	1532278	14.454	ng	99
78) Pyrene	19.45	202	1281422	9.976	ng	98
80) Butylbenzylphthalate	20.33	149	7414	3.191	ng	97
81) Benzo(a)anthracene	21.15	228	700477	5.597	ng	99
83) Chrysene	21.20	228	867587	7.349	ng	97
87) Indeno(1,2,3-cd)pyrene	25.70	276	707415	4.371	ng	95
88) Benzo(b)fluoranthene	22.74	252	1509297	10.583	ng	99
89) Benzo(k)fluoranthene	22.78	252	453700m	3.366	ng	
90) Benzo(a)pyrene	23.32	252	879276	6.823	ng	100
92) Benzo(a,h,i)perylene	26.40	276	721327	5.415	ng	99

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

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