

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091620\
 Data File : BP003306.D
 Acq On : 16 Sep 2020 17:35
 Operator : CG/JU
 Sample : L4018-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VNJ-229

Manual Integrations
 APPROVED

mohammad
 9/18/2020 11:28:36 AM

Quant Time: Sep 16 18:15:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 16 13:18:39 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.617	152	115816	20.00	ng	0.00
21) Naphthalene-d8	10.393	136	427743	20.00	ng	0.00
39) Acenaphthene-d10	14.263	164	236031	20.00	ng	0.00
64) Phenanthrene-d10	17.016	188	430394	20.00	ng	0.00
76) Chrysene-d12	21.110	240	404368	20.00	ng	0.00
86) Perylene-d12	23.322	264	545480	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.228	112	651069	98.16	ng	0.00
7) Phenol-d6	6.811	99	788472	89.19	ng	0.00
23) Nitrobenzene-d5	8.764	82	515048	70.74	ng	0.00
42) 2,4,6-Tribromophenol	15.757	330	279122	77.53	ng	0.00
45) 2-Fluorobiphenyl	12.881	172	1010016	62.58	ng	0.00
79) Terphenyl-d14	19.586	244	1044675	53.83	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.722	163	120076	6.37	ng	99
52) Acenaphthene	14.322	154	302661	22.33	ng	99
55) Dibenzofuran	14.663	168	293912	13.58	ng	96
58) Fluorene	15.316	166	516216	30.21	ng	99
71) Phenanthrene	17.057	178	2224483	95.00	ng	100
72) Anthracene	17.145	178	193717	8.39	ng	98
75) Fluoranthene	19.039	202	1673238	56.83	ng	99
78) Pyrene	19.386	202	1048790	41.59	ng	99
81) Benzo(a)anthracene	21.092	228	317127	12.13	ng	99
83) Chrysene	21.145	228	293735	11.70	ng	99
84) Bis(2-ethylhexyl)phtha...	21.033	149	43114	2.65	ng	100
88) Benzo(b)fluoranthene	22.657	252	267695	7.77	ng	98
89) Benzo(k)fluoranthene	22.698	252	83601m	2.53	ng	
90) Benzo(a)pyrene	23.222	252	132799	4.09	ng	99

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

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