

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
 Data File : BP003156.D
 Acq On : 28 Aug 2020 18:15
 Operator : CG/JU
 Sample : L3837-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 789UMR-TP11-0006-01

Manual Integrations
 APPROVED

mohammad
 8/31/2020 11:14:13 AM

Quant Time: Aug 28 19:00:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 24 15:32:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	205111	20.00	ng	-0.01
21) Naphthalene-d8	10.44	136	830851	20.00	ng	0.00
39) Acenaphthene-d10	14.30	164	528849	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	1173074	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	1361091	20.00	ng	-0.01
86) Perylene-d12	23.38	264	1506380	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	916461	79.00	ng	0.00
7) Phenol-d6	6.85	99	1103756	75.90	ng	0.00
23) Nitrobenzene-d5	8.80	82	644963	55.96	ng	0.00
42) 2,4,6-Tribromophenol	15.80	330	575365	80.44	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	1906176	52.78	ng	0.00
79) Terphenyl-d14	19.63	244	2958276	47.83	ng	-0.01
Target Compounds						
50) Dimethylphthalate	13.76	163	106470	2.640	ng	99
71) Phenanthrene	17.09	178	191050	3.006	ng	99
75) Fluoranthene	19.07	202	475041	6.466	ng	99
78) Pyrene	19.42	202	406053	4.597	ng	99
81) Benzo(a)anthracene	21.13	228	238797	2.801	ng	100
83) Chrysene	21.18	228	273707	3.353	ng	98
88) Benzo(b)fluoranthene	22.71	252	434972m	4.646	ng	
90) Benzo(a)pyrene	23.28	252	275845	3.176	ng	99
92) Benzo(g,h,i)perylene	26.34	276	189866	2.050	ng	98

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

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