

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007595.D
 Acq On : 22 Oct 2021 18:05
 Operator : CG/JU
 Sample : M4281-16
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY75

Quant Time: Oct 22 23:53:36 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 20 10:49:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	191182	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	774828	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	506578	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	1097633	20.000	ng/ul	0.00
79) Chrysene-d12	21.421	240	1148065	20.000	ng/ul	0.00
88) Perylene-d12	23.868	264	1196619	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	20843	4.306	ng/uL	0.00
4) Pyridine-d5	3.799	84	107783	8.134	ng/ul	0.00
7) Phenol-d5	7.105	99	106196	7.113	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	230240	25.855	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	277477	23.624	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	188824	16.419	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	163356	33.314	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	170844	37.301	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	333357	29.840	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	391610	25.922	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	1094220	31.221	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1259709	28.601	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	44869	7.979	ng/ul	0.00
60) Fluorene-d10	15.575	176	942493	32.135	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	29631	6.737	ng/ul	0.00
73) Anthracene-d10	17.433	188	1622549	34.808	ng/ul	0.00
81) Pyrene-d10	19.669	212	1990866	34.688	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	2021449	33.950	ng/ul	0.00

Target Compounds	Qvalue

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

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