

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007587.D
 Acq On : 22 Oct 2021 13:33
 Operator : CG/JU
 Sample : M4281-07
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY63

Quant Time: Oct 22 14:23:48 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	7.945	152	172930	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	702534	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.580	164	458012	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	990758	20.000	ng/ul	0.00
79) Chrysene-d12	21.421	240	1030708	20.000	ng/ul	0.00
88) Perylene-d12	23.868	264	1040943	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	23132	5.283	ng/uL	0.00
4) Pyridine-d5	3.793	84	139918	11.674	ng/ul	0.00
7) Phenol-d5	7.104	99	109376	8.099	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	274112	34.031	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	302561	28.479	ng/ul	0.00
15) 4-Methylphenol-d8	8.651	113	196393	18.879	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	179497	40.373	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	184278	44.374	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	344042	33.966	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	448424	32.738	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	1217170	38.412	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1383632	34.745	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	39700	7.808	ng/ul	0.00
60) Fluorene-d10	15.574	176	1053804	39.740	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	139503	35.140	ng/ul	0.00
73) Anthracene-d10	17.433	188	1816893	43.181	ng/ul	0.00
81) Pyrene-d10	19.668	212	2196365	42.626	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.715	264	2170381	41.903	ng/ul	0.00

Target Compounds Qvalue

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

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