

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007571.D
 Acq On : 22 Oct 2021 02:34
 Operator : CG/JU
 Sample : M4217-04
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGC2

Quant Time: Oct 22 03:06:19 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	220220	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	895329	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	580642	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.334	188	1245643	20.000	ng/ul	0.00
79) Chrysene-d12	21.422	240	1284612	20.000	ng/ul	0.00
88) Perylene-d12	23.868	264	1310143	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	7582	1.360	ng/uL	0.00
4) Pyridine-d5	3.799	84	43480	2.849	ng/ul	0.00
7) Phenol-d5	7.105	99	133637	7.770	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	79319	7.733	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	106668	7.884	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	101449	7.658	ng/ul	0.00
21) Nitrobenzene-d5	9.105	128	45597	8.047	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	44466	8.402	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	98986	7.668	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	124881	7.154	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	323170	8.045	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	373959	7.407	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	31853	4.942	ng/ul	-0.01
60) Fluorene-d10	15.575	176	286523	8.523	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	18315	3.669	ng/ul	0.00
73) Anthracene-d10	17.434	188	447450	8.458	ng/ul	0.00
81) Pyrene-d10	19.663	212	578046	9.001	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	531358	8.151	ng/ul	-0.01

Target Compounds Qvalue

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

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