

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072522\
 Data File : BP011084.D
 Acq On : 25 Jul 2022 10:25
 Operator : CG/JU
 Sample : PB146368BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK368

Quant Time: Jul 25 11:52:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.669	152	467416	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.463	136	2153945	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.334	164	1366751	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	2709222	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.216	240	1958243	20.000	ng/u1	#-0.01
88) Perylene-d12	23.563	264	1649673	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	77077	5.945	ng/uL	-0.02
4) Pyridine-d5	3.575	84	1028456	30.689	ng/u1	-0.02
7) Phenol-d5	6.846	99	1214440	31.652	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.011	67	817484	33.134	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.199	132	1036785	33.441	ng/u1	-0.01
15) 4-Methylphenol-d8	8.387	113	1044504	34.005	ng/u1	-0.01
21) Nitrobenzene-d5	8.834	128	535051	34.086	ng/u1	0.00
24) 2-Nitrophenol-d4	9.552	143	572316	36.847	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	928021	32.671	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.610	131	1518942	32.437	ng/u1	0.00
46) Dimethylphthalate-d6	13.751	166	3368045	36.056	ng/u1	-0.01
49) Acenaphthylene-d8	14.022	160	3935971	35.712	ng/u1	0.00
54) 4-Nitrophenol-d4	14.551	143	493826	26.121	ng/u1	0.00
60) Fluorene-d10	15.334	176	2738123	34.485	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	431280	32.171	ng/u1	0.00
73) Anthracene-d10	17.198	188	4050321	35.167	ng/u1	0.00
81) Pyrene-d10	19.457	212	4188131	40.722	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.410	264	2773065	34.999	ng/u1	-0.02

Target Compounds Qvalue

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

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