

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082322\
 Data File : BP011536.D
 Acq On : 23 Aug 2022 10:01
 Operator : CG/JU
 Sample : PB146666BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK666

Quant Time: Aug 23 10:32:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 01:58:37 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.552	152	66659	20.000 ng/u1	0.00
20) Naphthalene-d8	10.334	136	278970	20.000 ng/u1	-0.02
38) Acenaphthene-d10	14.222	164	161733	20.000 ng/u1	-0.01
64) Phenanthrene-d10	16.992	188	329433	20.000 ng/u1	-0.01
79) Chrysene-d12	21.122	240	305880	20.000 ng/u1	-0.01
88) Perylene-d12	23.410	264	318387	20.000 ng/u1	-0.02
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.064	96	10073	4.689 ng/uL	0.00
4) Pyridine-d5	3.475	84	128475	23.361 ng/u1	0.00
7) Phenol-d5	6.734	99	152608	24.913 ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.899	67	101447	25.491 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.081	132	124986	27.741 ng/u1	-0.01
15) 4-Methylphenol-d8	8.269	113	121963	24.875 ng/u1	-0.01
21) Nitrobenzene-d5	8.710	128	57354	28.557 ng/u1	-0.02
24) 2-Nitrophenol-d4	9.428	143	60735	29.714 ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.963	165	105291	27.907 ng/u1	-0.01
31) 4-Chloroaniline-d4	10.487	131	164192	26.143 ng/u1	-0.01
46) Dimethylphthalate-d6	13.645	166	366619	32.050 ng/u1	-0.01
49) Acenaphthylene-d8	13.910	160	383216	30.142 ng/u1	-0.01
54) 4-Nitrophenol-d4	14.451	143	54071	24.534 ng/u1	-0.01
60) Fluorene-d10	15.228	176	289603	30.093 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.363	200	45128	25.351 ng/u1	0.00
73) Anthracene-d10	17.092	188	476174	32.637 ng/u1	-0.01
81) Pyrene-d10	19.357	212	536323	34.655 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.263	264	522100	33.711 ng/u1	-0.02

Target Compounds Qvalue

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

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