

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012850.D
 Acq On : 29 Nov 2022 14:04
 Operator : CG/JU
 Sample : PB149236BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 29 21:58:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Nov 27 22:55:08 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	562145	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2578655	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	1739194	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.404	188	3852836	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	3714284	20.000	ng/ul	-0.01
88) Perylene-d12	23.998	264	3371692	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	69114	5.268	ng/uL	0.00
4) Pyridine-d5	3.817	84	1866559	47.546	ng/ul	0.00
7) Phenol-d5	7.164	99	1209924	26.193	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	781078	27.441	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	955289	27.140	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	971421	26.101	ng/ul	-0.01
21) Nitrobenzene-d5	9.175	128	455674	28.871	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.905	143	495359	29.559	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	970242	24.813	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	604950	11.298	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	3291107	27.492	ng/ul	0.00
49) Acenaphthylene-d8	14.340	160	3652105	26.146	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	488514	22.889	ng/ul	0.00
60) Fluorene-d10	15.640	176	2928106	27.929	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	395290	21.466	ng/ul	0.00
73) Anthracene-d10	17.504	188	4762736	28.082	ng/ul	0.00
81) Pyrene-d10	19.728	212	5548830	25.710	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	4550708	27.478	ng/ul	-0.01

Target Compounds Qvalue

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

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