

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012857.D
 Acq On : 29 Nov 2022 18:03
 Operator : CG/JU
 Sample : PB149233BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 29 22:58:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	519685	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2350517	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	1627779	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	3607407	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	3611692	20.000	ng/ul	0.00
88) Perylene-d12	23.998	264	3446680	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	72014	5.938	ng/uL	0.00
4) Pyridine-d5	3.822	84	959865	26.448	ng/ul	0.00
7) Phenol-d5	7.163	99	1194376	27.969	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	759314	28.856	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	961991	29.564	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	972647	28.269	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	464163	32.264	ng/ul	0.00
24) 2-Nitrophenol-d4	9.904	143	511407	33.478	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	1002462	28.125	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	1475052	30.221	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	3388799	30.246	ng/ul	0.00
49) Acenaphthylene-d8	14.339	160	3785116	28.953	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	524317	26.248	ng/ul	0.00
60) Fluorene-d10	15.639	176	3003512	30.609	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	429187	24.893	ng/ul	0.00
73) Anthracene-d10	17.498	188	4920281	30.985	ng/ul	0.00
81) Pyrene-d10	19.727	212	5740155	27.352	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	5130835	30.307	ng/ul	0.00

Target Compounds Qvalue

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

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