

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012843.D
 Acq On : 29 Nov 2022 10:05
 Operator : CG/JU
 Sample : PB149186BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 29 21:55:58 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	8.016	152	468124	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	2087302	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	1394601	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	3064406	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	2834157	20.000	ng/ul	0.00
88) Perylene-d12	24.004	264	2655642	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	57804	5.291	ng/uL	0.00
4) Pyridine-d5	3.817	84	766601	23.449	ng/ul	0.00
7) Phenol-d5	7.164	99	948783	24.665	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	629613	26.562	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	775227	26.448	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	771433	24.890	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	358864	28.090	ng/ul	0.00
24) 2-Nitrophenol-d4	9.911	143	390563	28.792	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	776223	24.524	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	1109583	25.600	ng/ul	0.00
46) Dimethylphthalate-d6	14.057	166	2703277	28.161	ng/ul	0.00
49) Acenaphthylene-d8	14.346	160	2925961	26.124	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	348163	20.343	ng/ul	0.00
60) Fluorene-d10	15.645	176	2358560	28.056	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	278893	19.042	ng/ul	0.00
73) Anthracene-d10	17.504	188	3881903	28.778	ng/ul	0.00
81) Pyrene-d10	19.733	212	4515199	27.418	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.839	264	3679319	28.206	ng/ul	0.00

Target Compounds	Qvalue

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

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