

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070623\
 Data File : BP016069.D
 Acq On : 07 Jul 2023 14:09
 Operator : MA/JU
 Sample : PB153637BL
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 07 20:27:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 06 23:18:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	211306	20.000	ng/u1	0.00
20) Naphthalene-d8	10.857	136	920915	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.681	164	609822	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	1400231	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1400833	20.000	ng/u1	0.02
88) Perylene-d12	24.092	264	1451636	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	41296	7.031	ng/uL	0.00
4) Pyridine-d5	3.799	84	480424	32.450	ng/u1	0.00
7) Phenol-d5	7.222	99	620927	34.344	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.358	67	437892	39.148	ng/u1	0.00
11) 2-Chlorophenol-d4	7.564	132	531906	38.974	ng/u1	0.00
15) 4-Methylphenol-d8	8.775	113	546878	38.290	ng/u1	0.00
21) Nitrobenzene-d5	9.228	128	265511	37.726	ng/u1	0.01
24) 2-Nitrophenol-d4	9.963	143	299900	36.510	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.546	165	444576	30.372	ng/u1	0.04
31) 4-Chloroaniline-d4	11.052	131	673067	35.795	ng/u1	0.03
46) Dimethylphthalate-d6	14.098	166	1917379	40.679	ng/u1	0.01
49) Acenaphthylene-d8	14.375	160	2068455	39.038	ng/u1	0.00
54) 4-Nitrophenol-d4	15.010	143	138891	22.329	ng/u1	0.05
60) Fluorene-d10	15.687	176	1595822	41.118	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.869	200	227198	26.384	ng/u1	0.02
73) Anthracene-d10	17.551	188	2553148	39.918	ng/u1	0.00
81) Pyrene-d10	19.781	212	3238217	35.401	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.916	264	2953931	40.339	ng/u1	0.00

Target Compounds	Qvalue

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

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