

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090623\
 Data File : BP016967.D
 Acq On : 06 Sep 2023 22:29
 Operator : MA/JU
 Sample : 04166-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH9G3

Quant Time: Sep 06 23:18:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	545987	20.000	ng/u1	0.00
20) Naphthalene-d8	10.810	136	2369324	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.639	164	1462830	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	3153830	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	2768780	20.000	ng/u1	0.00
88) Perylene-d12	24.051	264	2837575	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	57908	4.012	ng/uL	0.00
4) Pyridine-d5	3.763	84	380851	9.234	ng/u1	0.00
7) Phenol-d5	7.128	99	366917	7.356	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.322	67	1058362	33.921	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	962968	26.674	ng/u1	0.00
15) 4-Methylphenol-d8	8.687	113	651079	17.133	ng/u1	0.00
21) Nitrobenzene-d5	9.181	128	641547	34.902	ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	686739	34.465	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	1086636	29.539	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	570559	10.750	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	3918468	35.986	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	4382490	34.874	ng/u1	0.00
54) 4-Nitrophenol-d4	14.851	143	151130	7.105	ng/u1	0.00
60) Fluorene-d10	15.639	176	3349390	35.926	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	604891	31.871	ng/u1	0.00
73) Anthracene-d10	17.510	188	5398153	37.271	ng/u1	0.00
81) Pyrene-d10	19.745	212	6168169	36.575	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.886	264	5517436	38.327	ng/u1	0.00

Target Compounds Qvalue

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

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