

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072921\
 Data File : BP006424.D
 Acq On : 29 Jul 2021 17:07
 Operator : CG/JU
 Sample : PB137967BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK967

Quant Time: Jul 29 17:45:17 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 29 16:44:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	322222	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	1444560	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.410	164	852395	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.157	188	1654714	20.000	ng/u1	0.00
79) Chrysene-d12	21.257	240	1455555	20.000	ng/u1	0.00
88) Perylene-d12	23.603	264	1259027	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	67937	7.870	ng/uL	0.00
4) Pyridine-d5	3.699	84	981850	38.314	ng/u1	0.00
7) Phenol-d5	6.951	99	1154582	38.152	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	749065	39.781	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	899460	40.006	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	933170	39.211	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	454229	42.230	ng/u1	0.00
24) 2-Nitrophenol-d4	9.651	143	473273	46.082	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	768204	38.171	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	1246417	37.619	ng/u1	0.00
46) Dimethylphthalate-d6	13.828	166	2510785	41.386	ng/u1	0.00
49) Acenaphthylene-d8	14.098	160	3040074	40.787	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	443019	36.778	ng/u1	0.00
60) Fluorene-d10	15.404	176	2033513	39.873	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.522	200	302503	32.793	ng/u1	0.00
73) Anthracene-d10	17.263	188	3168967	42.425	ng/u1	0.00
81) Pyrene-d10	19.504	212	3422010	45.708	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.451	264	2730174	42.672	ng/u1	0.00

Target Compounds Qvalue

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

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